

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\
 Data File : BM029952.D
 Acq On : 11 May 2021 20:49
 Operator : CG/JU
 Sample : M2177-23
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG594

Manual Integrations
 APPROVED

mohammad
 5/12/2021 2:06:53 PM

Quant Time: May 12 01:26:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 11 15:21:38 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	123259	20.00	ng/ul	0.00
20) Naphthalene-d8	10.32	136	526622	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.19	164	320581	20.00	ng/ul	-0.01
64) Phenanthrene-d10	16.94	188	610185	20.00	ng/ul	0.00
79) Chrysene-d12	21.13	240	576982	20.00	ng/ul	0.00
88) Perylene-d12	23.29	264	590299	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	6789m	2.20	ng/uL	0.00
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul	
7) Phenol-d5	6.73	99	165293	14.42	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.89	67	115891	16.30	ng/ul	0.00
11) 2-Chlorophenol-d4	7.08	132	140610	15.89	ng/ul	0.00
15) 4-Methylphenol-d8	8.26	113	141477	14.92	ng/ul	0.00
21) Nitrobenzene-d5	8.70	128	61756	17.02	ng/ul	0.00
24) 2-Nitrophenol-d4	9.42	143	68870	16.81	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.96	165	134013	16.32	ng/ul	0.00
31) 4-Chloroaniline-d4	10.47	131	135898	10.87	ng/ul	0.00
46) Dimethylphthalate-d6	13.62	166	452174	18.20	ng/ul	0.00
49) Acenaphthylene-d8	13.89	160	565010	17.94	ng/ul	0.00
54) 4-Nitrophenol-d4	14.45	143	32891m	8.13	ng/ul	0.02
60) Fluorene-d10	15.19	176	411925	19.53	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	37958	9.58	ng/ul	0.00
73) Anthracene-d10	17.04	188	602032	19.59	ng/ul	0.00
81) Pyrene-d10	19.34	212	638461	21.57	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.15	264	669157	20.00	ng/ul	0.00

Target Compounds

					Ovalue	
37) 1-Methylnaphthalene	12.22	142	31727	1.510	ng/ul#	99
52) Acenaphthene	14.26	153	28227	1.311	ng/ul	94
61) Fluorene	15.25	166	31990	1.269	ng/ul#	93
72) Phenanthrene	16.99	178	337295	9.375	ng/ul	99
74) Anthracene	17.07	178	82597m	2.246	ng/ul	
80) Fluoranthene	19.00	202	348134	9.664	ng/ul	98
82) Pyrene	19.37	202	506674	13.365	ng/ul	99
85) Benzo(a)anthracene	21.12	228	205218	5.451	ng/ul#	91
86) Bis(2-ethylhexyl)phthalate	21.06	149	23966	1.032	ng/ul#	91
87) Chrysene	21.17	228	232867	6.211	ng/ul#	95
90) Benzo(b)fluoranthene	22.64	252	172156	4.496	ng/ul#	82
91) Benzo(k)fluoranthene	22.69	252	55253m	1.373	ng/ul	
93) Benzo(a)pyrene	23.19	252	161910	4.656	ng/ul#	89
94) Indeno(1,2,3-cd)pyrene	25.43	276	78843	1.873	ng/ul	95
96) Benzo(g,h,i)perylene	26.09	276	76585	2.246	ng/ul	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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