

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012920\
 Data File : BM024655.D
 Acq On : 29 Jan 2020 18:05
 Operator : CG/JU
 Sample : L1287-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 AOC-3-EXC-PIT-(0-5)-B

Manual Integrations
 APPROVED

mohammad
 1/30/2020 10:19:06 AM

Quant Time: Jan 30 09:07:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 23 14:57:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	422219	20.00	ng	0.00
21) Naphthalene-d8	10.20	136	1521150	20.00	ng	0.00
39) Acenaphthene-d10	14.10	164	873020	20.00	ng	0.00
64) Phenanthrene-d10	16.85	188	1484032	20.00	ng	0.00
76) Chrysene-d12	21.07	240	1348748	20.00	ng	0.01
87) Perylene-d12	23.19	264	1598306	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.09	112	1925831	86.74	ng	0.01
7) Phenol-d6	6.65	99	2435079	79.62	ng	0.00
23) Nitrobenzene-d5	8.59	82	1563217	50.40	ng	0.00
42) 2,4,6-Tribromophenol	15.60	330	997331	70.23	ng	0.00
45) 2-Fluorobiphenyl	12.71	172	3507322	54.62	ng	0.00
79) Terphenyl-d14	19.51	244	3871779	53.59	ng	0.00

Target Compounds

						Qvalue
10) Phenol	6.67	94	94103	3.087	ng	95
13) 1,4-Dichlorobenzene	7.47	146	67559	2.145	ng	95
31) Naphthalene	10.25	128	2413499	31.414	ng	100
37) 2-Methylnaphthalene	11.88	142	3767163	63.537	ng	99
38) 1-Methylnaphthalene	12.11	142	7549137	133.853	ng	100
46) 1,1'-Biphenyl	12.92	154	532392	8.058	ng	99
49) Acenaphthylene	13.82	152	3012680	37.731	ng	# 95
50) Dimethylphthalate	13.57	163	563839	7.987	ng	# 97
52) Acenaphthene	14.17	154	5454410	110.086	ng	99
55) Dibenzofuran	14.50	168	760670	9.414	ng	# 91
58) Fluorene	15.16	166	5177882	76.251	ng	98
71) Phenanthrene	16.90	178	15767354	196.202	ng	# 88
72) Anthracene	16.99	178	7006252	89.101	ng	99
73) Carbazole	17.26	167	364524	5.092	ng	# 79
74) Di-n-butylphthalate	17.85	149	247386	2.703	ng	# 74
75) Fluoranthene	18.93	202	10763181	107.309	ng	97
78) Pyrene	19.29	202	13789066	155.082	ng	93
80) Butylbenzylphthalate	20.22	149	293344	7.862	ng	# 94
81) Benzo(a)anthracene	21.05	228	6616859m	70.685	ng	
83) Chrysene	21.10	228	6171262	69.088	ng	97
84) Bis(2-ethylhexyl)phthalate	21.02	149	8654307	157.206	ng	# 96
86) Indeno(1,2,3-cd)pyrene	25.27	276	2473421	25.406	ng	97
88) Benzo(b)fluoranthene	22.57	252	5675130m	54.746	ng	
89) Benzo(k)fluoranthene	22.60	252	1632450m	16.133	ng	
90) Benzo(a)pyrene	23.10	252	5088076	53.986	ng	97
91) Dibenzo(a,h)anthracene	25.27	278	778261	8.700	ng	# 89
92) Benzo(g,h,i)perylene	25.90	276	2442013	30.065	ng	97

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

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