

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081916\
 Data File : BM007214.D
 Acq On : 18 Aug 2016 21:15
 Operator : UM/SJ
 Sample : H4445-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 E5N28

Manual Integrations
 APPROVED

sohil
 8/19/2016 6:41:17 PM

Quant Time: Aug 19 05:27:58 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 19 04:48:28 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	222192	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	898655	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	519050	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	1232314	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	1590592	20.00	ng/ul	0.00
83) Perylene-d12	23.63	264	1645694	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.32	96	18125	3.56	ng/uL	0.00
5) Phenol-d5	6.97	99	371807	20.66	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	212145	20.56	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	293699	20.49	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	298590	20.16	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	136570	20.92	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	150891	21.64	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.21	165	293483	20.09	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	353509	20.34	ng/ul	0.00
43) Dimethylphthalate-d6	13.84	166	924007	21.54	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	1109064	21.53	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	145476	18.92	ng/ul	0.00
57) Fluorene-d10	15.43	176	819853	21.87	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	117659	17.28	ng/ul	0.00
70) Anthracene-d10	17.27	188	1338981	23.45	ng/ul	0.00
76) Pyrene-d10	19.57	212	1667540	24.84	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.48	264	1846525	24.53	ng/ul	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
44) Dimethylphthalate	13.89	163	226772	5.29	ng/ul	98
47) Acenaphthylene	14.15	152	57538	1.08	ng/ul	96
69) Phenanthrene	17.22	178	80472	1.21	ng/ul	97
71) Anthracene	17.31	178	78217	1.15	ng/ul	99
74) Fluoranthene	19.23	202	493023	6.13	ng/ul	100
77) Pyrene	19.59	202	521023	5.92	ng/ul	99
80) Benzo(a)anthracene	21.34	228	348929	3.71	ng/ul	99
82) Chrysene	21.39	228	318888	3.71	ng/ul	98
85) Benzo(b)fluoranthene	22.94	252	602317	6.08	ng/ul	96
86) Benzo(k)fluoranthene	22.99	252	182341m	1.96	ng/ul	
88) Benzo(a)pyrene	23.53	252	325919	3.47	ng/ul	100
89) Indeno(1,2,3-cd)pyrene	25.91	276	232205	2.02	ng/ul	98
91) Benzo(g,h,i)perylene	26.61	276	187752	1.93	ng/ul	99

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

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