

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
 Data File : BP006983.D
 Acq On : 15 Sep 2021 19:38
 Operator : CG/JU
 Sample : M3608-03DL 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBKJ9DL

Manual Integrations
 APPROVED

mohammad

9/16/2021 11:44:09 AM

Quant Time: Sep 16 01:01:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 15 17:01:19 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	459409	20.000	ng/u1	0.00
20) Naphthalene-d8	10.722	136	2018438	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.546	164	1389402	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.292	188	2918411	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.375	240	2777823	20.000	ng/u1	# 0.00
88) Perylene-d12	23.792	264	2717423	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	2829m	0.245	ng/uL	0.00
4) Pyridine-d5	3.823	84	16767m	0.557	ng/u1	0.05
7) Phenol-d5	7.075	99	23829	0.706	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	60022	3.095	ng/u1	0.00
11) 2-Chlorophenol-d4	7.452	132	61585	2.259	ng/u1	0.00
15) 4-Methylphenol-d8	8.622	113	42603	1.607	ng/u1	0.00
21) Nitrobenzene-d5	9.081	128	34849	2.850	ng/u1	0.00
24) 2-Nitrophenol-d4	9.799	143	29193	2.425	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	69723	2.457	ng/u1	0.00
31) 4-Chloroaniline-d4	10.858	131	92759m	2.363	ng/u1	0.00
46) Dimethylphthalate-d6	13.957	166	327794	3.643	ng/u1	0.00
49) Acenaphthylene-d8	14.240	160	331153	2.906	ng/u1	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/u1	
60) Fluorene-d10	15.534	176	283862	3.574	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.657	200	76623	5.935	ng/u1	0.00
73) Anthracene-d10	17.392	188	462907	3.831	ng/u1	0.00
81) Pyrene-d10	19.622	212	512703	3.941	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.633	264	428492	3.269	ng/u1	-0.01
Target Compounds						
					Qvalue	
30) Naphthalene	10.769	128	2487997	25.411	ng/u1	100
36) 2-Methylnaphthalene	12.375	142	251171	3.826	ng/u1	98
37) 1-Methylnaphthalene	12.593	142	975398	14.566	ng/u1	99
43) 1,1'-Biphenyl	13.387	154	428471	4.429	ng/u1#	96
52) Acenaphthene	14.610	153	1306498	16.685	ng/u1	97
56) Dibenzofuran	14.940	168	1151474	10.191	ng/u1	87
58) 2,3,4,6-Tetrachlorophenol	15.169	232	27858	1.347	ng/u1#	77
61) Fluorene	15.593	166	691861	7.682	ng/u1	98
71) Pentachlorophenol	16.940	266	441026	24.066	ng/u1	99
72) Phenanthrene	17.334	178	1207972	8.309	ng/u1	98
77) Carbazole	17.692	167	1948440	15.171	ng/u1	97

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

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