

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007666.D
 Acq On : 26 Oct 2021 05:33
 Operator : CG/JU
 Sample : M4291-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P028-TW001-01

Quant Time: Oct 26 06:29:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 25 16:57:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	191116	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	772877	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	519389	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1152238	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1257361	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	1321617	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	21832	5.189	ng/uL	0.00
4) Pyridine-d5	3.793	84	49368	3.880	ng/ul	0.00
7) Phenol-d5	7.099	99	110414	7.606	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	267985	33.251	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	298611	26.449	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	196954	17.327	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	175444	33.310	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	180856	31.439	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	331157	28.282	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.869	131	362974	23.444	ng/ul	-0.01
46) Dimethylphthalate-d6	13.981	166	1196527	34.010	ng/ul	-0.01
49) Acenaphthylene-d8	14.263	160	1390867	32.255	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	43942	6.814	ng/ul	-0.01
60) Fluorene-d10	15.569	176	1068311	36.019	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	175771	28.519	ng/ul	0.00
73) Anthracene-d10	17.428	188	1934188	40.847	ng/ul	0.00
81) Pyrene-d10	19.663	212	2363876	40.381	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	2450132	37.739	ng/ul	0.00

Target Compounds Qvalue

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

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