

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007672.D
 Acq On : 26 Oct 2021 09:31
 Operator : CG/JU
 Sample : PB140137BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK137

Quant Time: Oct 26 10:55:03 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	204296	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	825433	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	543035	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.328	188	1173115	20.000	ng/u1	0.00
79) Chrysene-d12	21.416	240	1282454	20.000	ng/u1	0.00
88) Perylene-d12	23.863	264	1363747	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	29002	6.449	ng/uL	0.00
4) Pyridine-d5	3.781	84	429140	31.554	ng/u1	0.00
7) Phenol-d5	7.099	99	509485	32.833	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	299743	34.792	ng/u1	0.00
11) 2-Chlorophenol-d4	7.464	132	417305	34.577	ng/u1	-0.01
15) 4-Methylphenol-d8	8.640	113	404845	33.319	ng/u1	-0.01
21) Nitrobenzene-d5	9.093	128	195031	34.672	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	213773	34.795	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.358	165	400577	32.032	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.869	131	561019	33.929	ng/u1	-0.01
46) Dimethylphthalate-d6	13.981	166	1317469	35.817	ng/u1	-0.01
49) Acenaphthylene-d8	14.263	160	1527533	33.882	ng/u1	0.00
54) 4-Nitrophenol-d4	14.763	143	216741	32.145	ng/u1	-0.01
60) Fluorene-d10	15.569	176	1099562	35.458	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	168256	26.814	ng/u1	0.00
73) Anthracene-d10	17.428	188	1777720	36.874	ng/u1	0.00
81) Pyrene-d10	19.663	212	2176725	36.457	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.704	264	2262214	33.768	ng/u1	0.00

Target Compounds Qvalue

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

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