

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007684.D
 Acq On : 26 Oct 2021 18:27
 Operator : CG/JU
 Sample : PB140178BL
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK178

Quant Time: Oct 27 00:54:47 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	151221	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	618304	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.569	164	403455	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.328	188	878434	20.000	ng/u1	0.00
79) Chrysene-d12	21.416	240	944756	20.000	ng/u1	0.00
88) Perylene-d12	23.863	264	971694	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	25041	7.522	ng/uL	0.00
4) Pyridine-d5	3.781	84	365862	36.343	ng/u1	0.00
7) Phenol-d5	7.099	99	427368	37.208	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	250018	39.206	ng/u1	0.00
11) 2-Chlorophenol-d4	7.464	132	342580	38.348	ng/u1	-0.01
15) 4-Methylphenol-d8	8.640	113	336334	37.396	ng/u1	-0.01
21) Nitrobenzene-d5	9.093	128	159838	37.934	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	172557	37.496	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.358	165	331636	35.403	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.869	131	467278	37.727	ng/u1	-0.01
46) Dimethylphthalate-d6	13.981	166	1083297	39.639	ng/u1	-0.01
49) Acenaphthylene-d8	14.263	160	1250186	37.323	ng/u1	0.00
54) 4-Nitrophenol-d4	14.763	143	172620	34.458	ng/u1	-0.01
60) Fluorene-d10	15.569	176	911983	39.584	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	130375	27.747	ng/u1	0.00
73) Anthracene-d10	17.428	188	1464865	40.578	ng/u1	0.00
81) Pyrene-d10	19.663	212	1770545	40.254	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.704	264	1786628	37.429	ng/u1	0.00

Target Compounds Qvalue

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

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