

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102821\
 Data File : BP007746.D
 Acq On : 28 Oct 2021 10:59
 Operator : CG/JU
 Sample : PB140114BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK114

Quant Time: Oct 28 11:32:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 27 15:37:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	197309	20.000	ng/u1	0.00
20) Naphthalene-d8	10.734	136	834532	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.569	164	582089	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.322	188	1258969	20.000	ng/u1	0.00
79) Chrysene-d12	21.416	240	1319671	20.000	ng/u1	0.00
88) Perylene-d12	23.857	264	1375034	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	24411	5.620	ng/uL	0.00
4) Pyridine-d5	3.781	84	355547	27.068	ng/u1	0.00
7) Phenol-d5	7.093	99	442247	29.509	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	244636	29.401	ng/u1	0.00
11) 2-Chlorophenol-d4	7.464	132	366968	31.483	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	362943	30.928	ng/u1	-0.01
21) Nitrobenzene-d5	9.087	128	178070	31.311	ng/u1	0.00
24) 2-Nitrophenol-d4	9.810	143	203852	32.819	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	366586	28.995	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	509820	30.496	ng/u1	0.00
46) Dimethylphthalate-d6	13.975	166	1201199	30.465	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	1418174	29.346	ng/u1	0.00
54) 4-Nitrophenol-d4	14.757	143	199365	27.584	ng/u1	0.00
60) Fluorene-d10	15.563	176	1020395	30.698	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	163439	24.270	ng/u1	0.00
73) Anthracene-d10	17.422	188	1689718	32.659	ng/u1	0.00
81) Pyrene-d10	19.657	212	1993247	32.442	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.704	264	2054032	30.409	ng/u1	0.00

Target Compounds	Qvalue

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

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