

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007701.D
 Acq On : 27 Oct 2021 04:39
 Operator : CG/JU
 Sample : PB140187BL
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK187

Quant Time: Oct 27 05:55:05 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	171627	20.000	ng/u1	0.00
20) Naphthalene-d8	10.734	136	701203	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.569	164	460583	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.322	188	1009815	20.000	ng/u1	-0.01
79) Chrysene-d12	21.416	240	1094925	20.000	ng/u1	0.00
88) Perylene-d12	23.857	264	1150255	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	26690	7.064	ng/uL	0.00
4) Pyridine-d5	3.781	84	384137	33.621	ng/u1	0.00
7) Phenol-d5	7.093	99	454316	34.851	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.263	67	261116	36.078	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	367488	36.246	ng/u1	-0.01
15) 4-Methylphenol-d8	8.640	113	355494	34.826	ng/u1	-0.01
21) Nitrobenzene-d5	9.093	128	169810	35.536	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	179047	34.306	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	346115	32.581	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.869	131	490353	34.909	ng/u1	-0.01
46) Dimethylphthalate-d6	13.981	166	1145378	36.712	ng/u1	-0.01
49) Acenaphthylene-d8	14.263	160	1320051	34.521	ng/u1	0.00
54) 4-Nitrophenol-d4	14.757	143	181150	31.676	ng/u1	-0.02
60) Fluorene-d10	15.563	176	965849	36.722	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.681	200	137671	25.487	ng/u1	0.00
73) Anthracene-d10	17.422	188	1560640	37.606	ng/u1	-0.01
81) Pyrene-d10	19.657	212	1918570	37.637	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.704	264	1977606	34.998	ng/u1	0.00

Target Compounds	Qvalue

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

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