

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
 Data File : BP008532.D
 Acq On : 23 Dec 2021 19:24
 Operator : CG/JU
 Sample : PB141514BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK514

Quant Time: Dec 24 04:12:53 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 23 14:48:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	433698	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	1698842	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.533	164	1075755	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	2024751	20.000	ng/ul	0.00
79) Chrysene-d12	21.362	240	1786429	20.000	ng/ul	0.00
88) Perylene-d12	23.792	264	1979000	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	68824	7.153	ng/uL	0.00
4) Pyridine-d5	3.757	84	889468	33.033	ng/ul	0.00
7) Phenol-d5	7.057	99	1142772	33.030	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	641784	32.659	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	953184	34.632	ng/ul	0.00
15) 4-Methylphenol-d8	8.604	113	901958	32.497	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	447908	37.260	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	426511	38.695	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	908207	33.567	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	1292384	34.603	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	2726486	34.303	ng/ul	0.00
49) Acenaphthylene-d8	14.227	160	3496892	35.185	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	364263	28.911	ng/ul	0.00
60) Fluorene-d10	15.527	176	2346361	34.001	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	212249	23.814	ng/ul	0.00
73) Anthracene-d10	17.380	188	3506066	36.541	ng/ul	0.00
81) Pyrene-d10	19.610	212	3748464	35.018	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	3752753	36.844	ng/ul	0.00

Target Compounds Qvalue

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

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