

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018281.D
 Acq On : 22 Nov 2023 18:38
 Operator : MA/JU
 Sample : PB156975BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK975

Quant Time: Nov 22 23:19:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 22 14:09:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	381745	20.000	ng/u1	0.00
20) Naphthalene-d8	10.681	136	1639794	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.510	164	1109844	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	2380854	20.000	ng/u1	0.00
79) Chrysene-d12	21.363	240	1596731	20.000	ng/u1	0.00
88) Perylene-d12	23.780	264	1474495	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	52572	5.158	ng/uL	0.00
4) Pyridine-d5	3.687	84	746368	26.647	ng/u1	0.00
7) Phenol-d5	7.034	99	902660	25.873	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	565350	27.099	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	735308	25.103	ng/u1	0.00
15) 4-Methylphenol-d8	8.581	113	737612	26.198	ng/u1	0.00
21) Nitrobenzene-d5	9.040	128	355964	26.726	ng/u1	0.00
24) 2-Nitrophenol-d4	9.757	143	335890	27.307	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	716777	24.153	ng/u1	0.00
31) 4-Chloroaniline-d4	10.816	131	1144569	28.596	ng/u1	0.00
46) Dimethylphthalate-d6	13.928	166	2700221	27.343	ng/u1	0.00
49) Acenaphthylene-d8	14.204	160	2937857	26.109	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	323782	23.750	ng/u1	0.00
60) Fluorene-d10	15.504	176	2213321	26.796	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	190466	18.793	ng/u1	0.00
73) Anthracene-d10	17.357	188	3381368	26.572	ng/u1	0.00
81) Pyrene-d10	19.598	212	3498452	27.529	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.627	264	2275654	26.621	ng/u1	0.00

Target Compounds Qvalue

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

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