

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018262.D
 Acq On : 18 Nov 2023 15:30
 Operator : MA/JU
 Sample : PB157223BL
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK223

Quant Time: Nov 19 02:44:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	80080	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.593	136	307896	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.457	164	176482	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.239	188	408853	20.000	ng/u1	0.00
79) Chrysene-d12	21.380	240	379922	20.000	ng/u1	0.00
88) Perylene-d12	23.857	264	413266	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	14269	6.286	ng/uL	0.00
4) Pyridine-d5	3.587	84	158871	26.841	ng/u1	0.00
7) Phenol-d5	6.946	99	211027	27.296	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	139239	30.647	ng/u1	0.00
11) 2-Chlorophenol-d4	7.299	132	162156	26.486	ng/u1	0.00
15) 4-Methylphenol-d8	8.505	113	164081	25.900	ng/u1	0.00
21) Nitrobenzene-d5	8.981	128	83459	30.086	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.699	143	91598	29.147	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	144594	25.815	ng/u1	0.00
31) 4-Chloroaniline-d4	10.781	131	201924	26.288	ng/u1	0.00
46) Dimethylphthalate-d6	13.881	166	503449	30.918	ng/u1	0.00
49) Acenaphthylene-d8	14.157	160	530418	29.907	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	53354	21.927	ng/u1	0.00
60) Fluorene-d10	15.463	176	422405	31.420	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	70508	23.855	ng/u1	0.00
73) Anthracene-d10	17.345	188	701357	31.718	ng/u1	0.00
81) Pyrene-d10	19.598	212	829666	32.693	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.692	264	793188	33.008	ng/u1	0.00

Target Compounds Qvalue

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

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