

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018040.D
 Acq On : 10 Nov 2023 23:58
 Operator : MA/JU
 Sample : PB156684BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK684

Quant Time: Nov 11 00:47:21 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	43137	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	166384	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	97635	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.304	188	216046	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	204683	20.000	ng/u1	0.00
88) Perylene-d12	23.939	264	243158	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	9280	7.589	ng/uL	0.00
4) Pyridine-d5	3.664	84	107882	33.836	ng/u1	0.00
7) Phenol-d5	7.016	99	128374	30.825	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	80274	32.800	ng/u1	0.00
11) 2-Chlorophenol-d4	7.375	132	107756	32.674	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	102532	30.045	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	50530	33.708	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	56153	33.065	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	88213	29.144	ng/u1	0.00
31) 4-Chloroaniline-d4	10.851	131	128026	30.843	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	315595	35.033	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	333981	34.038	ng/u1	0.00
54) 4-Nitrophenol-d4	14.798	143	32832	24.390	ng/u1	0.02
60) Fluorene-d10	15.528	176	258222	34.719	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	43466	27.830	ng/u1	0.00
73) Anthracene-d10	17.410	188	413528	35.391	ng/u1	0.00
81) Pyrene-d10	19.657	212	495326	36.229	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.768	264	522140	36.929	ng/u1	0.00

Target Compounds Qvalue

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

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