

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017139.D
 Acq On : 13 Sep 2023 09:22
 Operator : MA/JU
 Sample : PB155418BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK418

Quant Time: Sep 13 23:24:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:27:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	170962	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.769	136	739619	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.610	164	484250	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	1144840	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	1147835	20.000	ng/u1	0.00
88) Perylene-d12	24.015	264	1263579	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	28044	6.893	ng/uL	-0.01
4) Pyridine-d5	3.740	84	331770	28.507	ng/u1	-0.01
7) Phenol-d5	7.087	99	475068	34.174	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.281	67	300657	34.603	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.463	132	374949	33.810	ng/u1	-0.01
15) 4-Methylphenol-d8	8.652	113	382283	33.181	ng/u1	-0.02
21) Nitrobenzene-d5	9.140	128	180999	34.949	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.857	143	196709	35.578	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.381	165	365937	34.223	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.934	131	264607	16.418	ng/u1	-0.01
46) Dimethylphthalate-d6	14.028	166	1284386	37.272	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	1363444	34.869	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	200192	30.058	ng/u1	0.00
60) Fluorene-d10	15.610	176	1111837	37.315	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	170724	26.691	ng/u1	0.00
73) Anthracene-d10	17.486	188	1841727	37.863	ng/u1	0.00
81) Pyrene-d10	19.727	212	2279120	37.860	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.851	264	2219242	37.211	ng/u1	0.00

Target Compounds Qvalue

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

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