

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017282.D
 Acq On : 22 Sep 2023 10:04
 Operator : MA/JU
 Sample : PB155610BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK610

Quant Time: Sep 22 10:45:15 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	151514	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	599597	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	345129	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	781434	20.000	ng/u1	0.00
79) Chrysene-d12	21.469	240	790252	20.000	ng/u1	0.00
88) Perylene-d12	23.992	264	893490	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	28522	7.732	ng/uL	0.00
4) Pyridine-d5	3.740	84	394437	38.249	ng/u1	0.00
7) Phenol-d5	7.093	99	467763	37.590	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	280060	37.293	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	379963	38.498	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	358683	37.075	ng/u1	0.00
21) Nitrobenzene-d5	9.134	128	173781	40.090	ng/u1	0.00
24) 2-Nitrophenol-d4	9.846	143	190846	40.548	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	342331	38.497	ng/u1	0.00
31) 4-Chloroaniline-d4	10.922	131	484984	37.650	ng/u1	0.00
46) Dimethylphthalate-d6	14.010	166	996208	40.293	ng/u1	0.00
49) Acenaphthylene-d8	14.298	160	1093606	38.459	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	156563	36.348	ng/u1	0.00
60) Fluorene-d10	15.592	176	862387	40.516	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.804	200	217	0.049	ng/u1	0.08
73) Anthracene-d10	17.469	188	1459708	42.119	ng/u1	0.00
81) Pyrene-d10	19.710	212	1806835	42.280	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.821	264	1843550	42.854	ng/u1	-0.01

Target Compounds Qvalue

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

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