

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017065.D
 Acq On : 09 Sep 2023 23:14
 Operator : MA/JU
 Sample : PB155240BL
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK240

Quant Time: Sep 10 01:06:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 09 01:41:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	769781	20.000	ng/u1	0.00
20) Naphthalene-d8	10.798	136	3349317	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	2012194	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.392	188	4214300	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	3207670	20.000	ng/u1	0.00
88) Perylene-d12	24.021	264	3024685	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	150953	7.418	ng/uL	0.00
4) Pyridine-d5	3.758	84	2429177	41.775	ng/u1	0.00
7) Phenol-d5	7.122	99	2879737	40.948	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	1833743	41.686	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	2128045	41.809	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	2214063	41.324	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	1069748	41.169	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	1175416	41.730	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	1978357	38.044	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	2931932	39.079	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	6215990	41.500	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	6991679	40.447	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	1107767	37.862	ng/u1	0.00
60) Fluorene-d10	15.622	176	5222669	40.725	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	829422	32.704	ng/u1	0.00
73) Anthracene-d10	17.498	188	8173861	42.234	ng/u1	0.00
81) Pyrene-d10	19.733	212	8823381	45.160	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.862	264	6909771	45.030	ng/u1	0.00

Target Compounds	Qvalue

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

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