

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017026.D
 Acq On : 08 Sep 2023 23:42
 Operator : MA/JU
 Sample : PB155208BL
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK208

Quant Time: Sep 09 04:34:38 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	7.975	152	420077	20.000	ng/u1	0.00
20) Naphthalene-d8	10.798	136	1839364	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	1117756	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	2303648	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	1662895	20.000	ng/u1	0.00
88) Perylene-d12	24.027	264	1560131	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	93246	8.396	ng/uL	0.00
4) Pyridine-d5	3.758	84	955773	30.120	ng/u1	0.00
7) Phenol-d5	7.122	99	1557441	40.581	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	1036309	43.169	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	1171387	42.173	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	1221817	41.789	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	592422	41.515	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	646560	41.798	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	1101207	38.560	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	776789	18.853	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	3516163	42.260	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	3909827	40.718	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	600236	36.932	ng/u1	0.00
60) Fluorene-d10	15.628	176	2932105	41.160	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	423583	30.554	ng/u1	0.00
73) Anthracene-d10	17.498	188	4499448	42.531	ng/u1	0.00
81) Pyrene-d10	19.733	212	4686864	46.273	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.862	264	3478174	43.945	ng/u1	0.00

Target Compounds Qvalue

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

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