

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
 Data File : BP021815.D
 Acq On : 04 Sep 2024 11:56
 Operator : RC/JU
 Sample : PB163092BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK092

Quant Time: Sep 04 12:55:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 29 23:56:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	437085	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.557	136	1792515	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.416	164	1080998	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.216	188	2389944	20.000	ng/u1	-0.01
79) Chrysene-d12	21.663	240	2279862	20.000	ng/u1	0.00
88) Perylene-d12	25.074	264	2478359	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	61715	4.688	ng/uL	0.00
4) Pyridine-d5	3.687	84	799675	21.000	ng/u1	0.00
7) Phenol-d5	6.952	99	1017323	21.809	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.111	67	564413	22.234	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	686918	22.810	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	772824	21.441	ng/u1	-0.01
21) Nitrobenzene-d5	8.922	128	342633	23.758	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.640	143	366100	25.009	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.181	165	643095	22.293	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.687	131	907668	21.627	ng/u1	-0.01
46) Dimethylphthalate-d6	13.816	166	2054913	24.843	ng/u1	-0.02
49) Acenaphthylene-d8	14.104	160	2173779	23.407	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.616	143	349579	22.214	ng/u1	-0.01
60) Fluorene-d10	15.422	176	1698928	24.431	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.539	200	280407	21.332	ng/u1	-0.02
73) Anthracene-d10	17.322	188	2725499	24.029	ng/u1	-0.02
81) Pyrene-d10	19.686	212	3306816	25.373	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.839	264	3171706	24.000	ng/u1	0.00

Target Compounds Qvalue

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

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