

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
 Data File : BP022043.D
 Acq On : 26 Sep 2024 01:36
 Operator : RC/JU
 Sample : PB163537BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK537

Quant Time: Sep 26 03:21:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	87405	20.000	ng/u1	0.00
20) Naphthalene-d8	10.463	136	320592	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	247642	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.110	188	615784	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	723411	20.000	ng/u1	0.00
88) Perylene-d12	24.816	264	850175	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	16973	7.606	ng/uL	0.00
4) Pyridine-d5	3.634	84	232633	36.472	ng/u1	0.00
7) Phenol-d5	6.887	99	268187	37.510	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	170454	38.717	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	200836	38.824	ng/u1	0.00
15) 4-Methylphenol-d8	8.405	113	213452	37.676	ng/u1	0.00
21) Nitrobenzene-d5	8.840	128	99063	41.285	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	115539	42.467	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	234107	39.623	ng/u1	0.00
31) 4-Chloroaniline-d4	10.599	131	275714	39.110	ng/u1	0.00
46) Dimethylphthalate-d6	13.722	166	809338	42.278	ng/u1	0.00
49) Acenaphthylene-d8	14.004	160	815255	40.224	ng/u1	0.00
54) 4-Nitrophenol-d4	14.528	143	121284	37.040	ng/u1	0.00
60) Fluorene-d10	15.322	176	697111	41.168	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.440	200	129221	33.800	ng/u1	0.00
73) Anthracene-d10	17.216	188	1201224	41.786	ng/u1	0.00
81) Pyrene-d10	19.586	212	1581815	42.206	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.592	264	1804618	40.357	ng/u1	-0.02

Target Compounds Qvalue

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

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