

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021686.D
 Acq On : 28 Aug 2024 02:19
 Operator : RC/JU
 Sample : PB162994BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK994

Quant Time: Aug 28 02:53:27 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 27 05:11:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	356275	20.000	ng/u1	0.00
20) Naphthalene-d8	10.575	136	1516771	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	989795	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.240	188	2136769	20.000	ng/u1	-0.01
79) Chrysene-d12	21.692	240	2074592	20.000	ng/u1	0.00
88) Perylene-d12	25.110	264	2258488	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	57301	5.340	ng/uL	0.00
4) Pyridine-d5	3.687	84	793945	25.578	ng/u1	0.00
7) Phenol-d5	6.958	99	1017439	26.759	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	584915	28.268	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	666446	27.150	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	776492	26.429	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	341895	28.017	ng/u1	0.00
24) 2-Nitrophenol-d4	9.658	143	350121	28.266	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	647065	26.509	ng/u1	0.00
31) 4-Chloroaniline-d4	10.705	131	894621	25.192	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	2184790	28.847	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2394435	28.159	ng/u1	0.00
54) 4-Nitrophenol-d4	14.628	143	370537	25.716	ng/u1	0.00
60) Fluorene-d10	15.440	176	1814953	28.504	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.557	200	241524	20.551	ng/u1	-0.01
73) Anthracene-d10	17.340	188	2937624	28.968	ng/u1	-0.01
81) Pyrene-d10	19.710	212	3588652	30.260	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.863	264	3462950	28.755	ng/u1	-0.01

Target Compounds Qvalue

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

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