

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
 Data File : BP021716.D
 Acq On : 29 Aug 2024 10:24
 Operator : RC/JU
 Sample : PB162969BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK969

Quant Time: Aug 30 00:20:15 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	7.787	152	348970	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	1479496	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.434	164	955036	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.234	188	2021114	20.000	ng/u1	-0.02
79) Chrysene-d12	21.692	240	1930558	20.000	ng/u1	0.00
88) Perylene-d12	25.092	264	2120917	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	60327	5.739	ng/uL	-0.01
4) Pyridine-d5	3.681	84	803722	26.435	ng/u1	-0.01
7) Phenol-d5	6.952	99	1022583	27.458	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.117	67	584359	28.832	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.322	132	671177	27.915	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	774528	26.914	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	337144	28.324	ng/u1	0.00
24) 2-Nitrophenol-d4	9.652	143	343024	28.390	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	630762	26.492	ng/u1	0.00
31) 4-Chloroaniline-d4	10.705	131	887302	25.615	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	2079907	28.462	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2292578	27.943	ng/u1	0.00
54) 4-Nitrophenol-d4	14.628	143	340957	24.524	ng/u1	0.00
60) Fluorene-d10	15.440	176	1740920	28.337	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.563	200	218500	19.656	ng/u1	0.00
73) Anthracene-d10	17.340	188	2803409	29.227	ng/u1	-0.01
81) Pyrene-d10	19.704	212	3354541	30.397	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.845	264	3216907	28.444	ng/u1	-0.03

Target Compounds Qvalue

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

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