

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021657.D
 Acq On : 27 Aug 2024 04:46
 Operator : RC/JU
 Sample : PB162966BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK966

Quant Time: Aug 27 05:23:15 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	7.793	152	308714	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1234317	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.439	164	797799	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1751246	20.000	ng/u1	-0.02
79) Chrysene-d12	21.686	240	1778706	20.000	ng/u1	0.00
88) Perylene-d12	25.098	264	2037506	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	52088	5.602	ng/uL	0.00
4) Pyridine-d5	3.693	84	713139	26.515	ng/u1	0.00
7) Phenol-d5	6.958	99	874404	26.540	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	500987	27.942	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	583533	27.434	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	660545	25.946	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	286908	28.891	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	291892	28.957	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	540936	27.232	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	743273	25.719	ng/u1	0.00
46) Dimethylphthalate-d6	13.839	166	1814790	29.728	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	1984938	28.961	ng/u1	0.00
54) 4-Nitrophenol-d4	14.628	143	308564	26.568	ng/u1	0.00
60) Fluorene-d10	15.445	176	1548998	30.182	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	230111	23.890	ng/u1	-0.01
73) Anthracene-d10	17.339	188	2493212	29.998	ng/u1	-0.01
81) Pyrene-d10	19.704	212	3097026	30.459	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.868	264	3209191	29.538	ng/u1	0.00

Target Compounds	Qvalue

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

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