

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021660.D
 Acq On : 27 Aug 2024 06:53
 Operator : RC/JU
 Sample : P3711-02
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E2494

Quant Time: Aug 27 07:38:08 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	7.792	152	307283	20.000	ng/u1	0.00
20) Naphthalene-d8	10.575	136	1286143	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.433	164	857574	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1911388	20.000	ng/u1	-0.02
79) Chrysene-d12	21.692	240	1904057	20.000	ng/u1	0.00
88) Perylene-d12	25.103	264	2144864	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	43778	4.730	ng/uL	0.00
4) Pyridine-d5	3.693	84	155589	5.812	ng/u1	0.00
7) Phenol-d5	6.957	99	300578	9.166	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	532751	29.852	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	578222	27.311	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	507048	20.009	ng/u1	0.00
21) Nitrobenzene-d5	8.939	128	309893	29.948	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	323318	30.782	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	582440	28.140	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	706078	23.448	ng/u1	0.00
46) Dimethylphthalate-d6	13.833	166	2062583	31.433	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2195764	29.804	ng/u1	0.00
54) 4-Nitrophenol-d4	14.627	143	103790	8.314	ng/u1	0.00
60) Fluorene-d10	15.439	176	1721152	31.199	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.557	200	273053	25.973	ng/u1	-0.01
73) Anthracene-d10	17.339	188	2983972	32.895	ng/u1	-0.01
81) Pyrene-d10	19.715	212	3598645	33.063	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.868	264	3784996	33.094	ng/u1	0.00

Target Compounds	Qvalue

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

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