

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021649.D
 Acq On : 26 Aug 2024 22:22
 Operator : RC/JU
 Sample : P3711-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E24A0

Quant Time: Aug 26 23:22:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 23 17:51:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	384639	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	1652472	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.445	164	1111213	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	2463563	20.000	ng/u1	0.01
79) Chrysene-d12	21.692	240	2381799	20.000	ng/u1	-0.02
88) Perylene-d12	25.104	264	2624163	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	43352	3.742	ng/uL	0.00
4) Pyridine-d5	3.693	84	203245	6.065	ng/u1	0.00
7) Phenol-d5	6.963	99	311688	7.593	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	564030	25.249	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	616709	23.271	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	542792	17.112	ng/u1	0.00
21) Nitrobenzene-d5	8.940	128	352707	26.530	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	375112	27.796	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	666020	25.044	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	775152	20.035	ng/u1	0.00
46) Dimethylphthalate-d6	13.839	166	2364289	27.806	ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	2544489	26.654	ng/u1	0.00
54) 4-Nitrophenol-d4	14.639	143	119871	7.410	ng/u1	0.01
60) Fluorene-d10	15.457	176	2024157	28.316	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.569	200	351172	25.917	ng/u1	0.00
73) Anthracene-d10	17.351	188	3576945	30.594	ng/u1	0.00
81) Pyrene-d10	19.721	212	4354636	31.983	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.862	264	4435708	31.700	ng/u1	-0.04

Target Compounds Qvalue

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

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