

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022813.D
 Acq On : 01 Nov 2024 20:52
 Operator : RC/JU
 Sample : P4627-04
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CR2

Quant Time: Nov 04 04:49:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	53653	20.000	ng/u1	0.00
20) Naphthalene-d8	10.387	136	200039	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.257	164	156634	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.063	188	376354	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	447355	20.000	ng/u1	0.00
88) Perylene-d12	24.763	264	557305	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	6234	4.515	ng/uL	0.00
4) Pyridine-d5	3.599	84	33155	8.747	ng/u1	0.00
7) Phenol-d5	6.828	99	43742	9.685	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	81558	30.725	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.175	132	95510	29.803	ng/u1	0.00
15) 4-Methylphenol-d8	8.340	113	79332	22.011	ng/u1	0.00
21) Nitrobenzene-d5	8.775	128	53420	34.731	ng/u1	0.00
24) 2-Nitrophenol-d4	9.487	143	64712	36.340	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.028	165	130848	34.573	ng/u1	0.00
31) 4-Chloroaniline-d4	10.534	131	122601	27.415	ng/u1	0.00
46) Dimethylphthalate-d6	13.663	166	468616	37.655	ng/u1	0.00
49) Acenaphthylene-d8	13.945	160	476015	36.013	ng/u1	0.00
54) 4-Nitrophenol-d4	14.534	143	16780	8.037	ng/u1	0.02
60) Fluorene-d10	15.275	176	421870	39.083	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.422	200	73498	29.693	ng/u1	0.02
73) Anthracene-d10	17.163	188	773293	43.678	ng/u1	0.00
81) Pyrene-d10	19.557	212	1034910	43.746	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.545	264	1311970	44.081	ng/u1	0.02

Target Compounds	Qvalue

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

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