

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102824\
 Data File : BP022677.D
 Acq On : 28 Oct 2024 12:58
 Operator : RC/JU
 Sample : P4529-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 F7H11

Quant Time: Oct 28 14:09:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 26 05:55:04 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.645	152	70334	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.404	136	259853	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.257	164	207486	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.057	188	508573	20.000	ng/u1	-0.02
79) Chrysene-d12	21.504	240	595137	20.000	ng/u1	0.00
88) Perylene-d12	24.762	264	727733	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	8573	4.737	ng/uL	-0.02
4) Pyridine-d5	3.593	84	128317	25.825	ng/u1	-0.02
7) Phenol-d5	6.840	99	163686	27.647	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67	99182	28.503	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.193	132	120760	28.745	ng/u1	-0.01
15) 4-Methylphenol-d8	8.351	113	139987	29.628	ng/u1	-0.01
21) Nitrobenzene-d5	8.787	128	61817	30.939	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.498	143	76050	32.877	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.039	165	150048	30.520	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.539	131	173383	29.846	ng/u1	-0.02
46) Dimethylphthalate-d6	13.669	166	574186	34.830	ng/u1	-0.01
49) Acenaphthylene-d8	13.945	160	577227	32.967	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.498	143	69196	25.020	ng/u1	0.00
60) Fluorene-d10	15.269	176	500499	35.003	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.416	200	87584	26.185	ng/u1	0.00
73) Anthracene-d10	17.174	188	879286	36.753	ng/u1	0.00
81) Pyrene-d10	19.545	212	1088194	34.577	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.527	264	1393301	35.850	ng/u1	0.00

Target Compounds Qvalue

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

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