

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102824\
 Data File : BP022680.D
 Acq On : 28 Oct 2024 14:58
 Operator : RC/JU
 Sample : P4529-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 F7H12

Quant Time: Oct 28 15:24:48 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	7.646	152	84697	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.399	136	316412	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.257	164	248403	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.057	188	620095	20.000	ng/u1	-0.02
79) Chrysene-d12	21.498	240	669373	20.000	ng/u1	0.00
88) Perylene-d12	24.757	264	782143	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	12115	5.559	ng/uL	-0.01
4) Pyridine-d5	3.599	84	169971	28.407	ng/u1	-0.01
7) Phenol-d5	6.840	99	224133	31.437	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67	130815	31.218	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.187	132	163734	32.365	ng/u1	-0.02
15) 4-Methylphenol-d8	8.352	113	186883	32.846	ng/u1	-0.01
21) Nitrobenzene-d5	8.787	128	84898	34.896	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.499	143	102106	36.251	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.034	165	206363	34.472	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.546	131	230475	32.582	ng/u1	-0.01
46) Dimethylphthalate-d6	13.663	166	757776	38.395	ng/u1	-0.02
49) Acenaphthylene-d8	13.945	160	763277	36.412	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.510	143	92943	28.071	ng/u1	0.00
60) Fluorene-d10	15.263	176	646118	37.744	ng/u1	-0.03
65) 4,6-Dinitro-2-methylph...	15.404	200	114703	28.125	ng/u1	-0.01
73) Anthracene-d10	17.163	188	1098026	37.641	ng/u1	0.00
81) Pyrene-d10	19.545	212	1359822	38.416	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.527	264	1620259	38.790	ng/u1	0.00

Target Compounds	Qvalue

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

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