

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102824\
 Data File : BP022695.D
 Acq On : 29 Oct 2024 01:03
 Operator : RC/JU
 Sample : P4530-13
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 F7H40

Quant Time: Oct 29 01:42:02 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	78533	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.405	136	290062	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.263	164	228461	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.057	188	574139	20.000	ng/u1	-0.02
79) Chrysene-d12	21.486	240	681721	20.000	ng/u1	-0.01
88) Perylene-d12	24.739	264	841114	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	6588	3.260	ng/uL	-0.02
4) Pyridine-d5	3.593	84	80985	14.597	ng/u1	-0.02
7) Phenol-d5	6.840	99	110741	16.752	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67	65819	16.940	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.187	132	80998	17.268	ng/u1	-0.02
15) 4-Methylphenol-d8	8.352	113	90023	17.064	ng/u1	-0.01
21) Nitrobenzene-d5	8.787	128	40071	17.967	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.499	143	48644	18.839	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.046	165	95816	17.459	ng/u1	0.00
31) 4-Chloroaniline-d4	10.546	131	105954	16.339	ng/u1	-0.01
46) Dimethylphthalate-d6	13.669	166	354341	19.521	ng/u1	-0.01
49) Acenaphthylene-d8	13.951	160	359792	18.662	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.510	143	42230	13.868	ng/u1	0.00
60) Fluorene-d10	15.269	176	308536	19.597	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.428	200	49242	13.041	ng/u1	0.01
73) Anthracene-d10	17.169	188	521223	19.298	ng/u1	0.00
81) Pyrene-d10	19.534	212	672782	18.662	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.516	264	886222	19.729	ng/u1	-0.01

Target Compounds	Qvalue

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

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