

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121224\
 Data File : BP023397.D
 Acq On : 12 Dec 2024 15:50
 Operator : RC/JU
 Sample : P5153-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E28P2

Quant Time: Dec 13 01:40:50 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.540	152	93337	20.000	ng/u1	0.00
20) Naphthalene-d8	10.287	136	347296	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.157	164	255369	20.000	ng/u1	0.01
64) Phenanthrene-d10	16.951	188	586998	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	778294	20.000	ng/u1	0.00
88) Perylene-d12	24.569	264	884979	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	7281	2.872	ng/uL	0.00
4) Pyridine-d5	3.523	84	78597	12.056	ng/u1	0.00
7) Phenol-d5	6.752	99	115969	15.835	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.893	67	77258	16.721	ng/u1	0.00
11) 2-Chlorophenol-d4	7.093	132	86112	15.931	ng/u1	0.00
15) 4-Methylphenol-d8	8.264	113	97197m	16.444	ng/u1	0.00
21) Nitrobenzene-d5	8.687	128	41863	16.142	ng/u1	0.00
24) 2-Nitrophenol-d4	9.399	143	49858	16.681	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.952	165	103644m	16.430	ng/u1	0.02
31) 4-Chloroaniline-d4	10.446	131	106493	15.069	ng/u1	0.00
46) Dimethylphthalate-d6	13.563	166	358848	18.101	ng/u1	0.00
49) Acenaphthylene-d8	13.846	160	366244	17.515	ng/u1	0.01
54) 4-Nitrophenol-d4	14.475	143	43012m	17.705	ng/u1	0.06
60) Fluorene-d10	15.163	176	301074	17.266	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.334	200	33437	9.425	ng/u1	0.02
73) Anthracene-d10	17.063	188	495135	18.482	ng/u1	0.01
81) Pyrene-d10	19.439	212	697379	17.150	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.351	264	803711	16.965	ng/u1	0.00

Target Compounds	Qvalue

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

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