

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
 Data File : BP019963.D
 Acq On : 22 Mar 2024 19:58
 Operator : MA/JU
 Sample : P1813-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0LX8

Quant Time: Mar 22 23:03:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 22 22:40:53 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	96893	20.000	ng/u1	0.00
20) Naphthalene-d8	10.375	136	383496	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.239	164	232931	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.080	188	448770	20.000	ng/u1	-0.01
79) Chrysene-d12	21.545	240	402781	20.000	ng/u1	0.00
88) Perylene-d12	24.845	264	472761	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	13187	5.963	ng/uL	0.00
4) Pyridine-d5	3.652	84	70186	10.511	ng/u1	0.00
7) Phenol-d5	6.810	99	54934	6.722	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	176238	33.102	ng/u1	0.00
11) 2-Chlorophenol-d4	7.169	132	146938	25.038	ng/u1	0.01
15) 4-Methylphenol-d8	8.322	113	103646	16.394	ng/u1	0.00
21) Nitrobenzene-d5	8.775	128	97569	33.193	ng/u1	0.00
24) 2-Nitrophenol-d4	9.481	143	104103	31.654	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	177005	29.470	ng/u1	0.01
31) 4-Chloroaniline-d4	10.528	131	155607	19.873	ng/u1	0.00
46) Dimethylphthalate-d6	13.675	166	639604	39.164	ng/u1	0.00
49) Acenaphthylene-d8	13.934	160	710625	37.249	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.457	143	28	0.011	ng/u1	-0.04
60) Fluorene-d10	15.275	176	533594	38.355	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200	86141	32.067	ng/u1	-0.02
73) Anthracene-d10	17.186	188	809967	40.773	ng/u1	0.00
81) Pyrene-d10	19.574	212	941712	39.849	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.615	264	942318	41.489	ng/u1	-0.02

Target Compounds Qvalue

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

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