

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011925.D
 Acq On : 05 Oct 2022 12:13
 Operator : CG/JU
 Sample : N4890-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EY0D9

Quant Time: Oct 05 23:21:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 05 23:18:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	242710	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	969311	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	566655	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.287	188	1243115	20.000	ng/u1	0.00
79) Chrysene-d12	21.369	240	1415891	20.000	ng/u1	0.00
88) Perylene-d12	23.804	264	1462759	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	27050	4.915	ng/uL	0.00
4) Pyridine-d5	3.728	84	121638	7.546	ng/u1	0.00
7) Phenol-d5	7.046	99	143054	6.827	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	361738	31.146	ng/u1	0.00
11) 2-Chlorophenol-d4	7.411	132	424315	25.653	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	254526	14.695	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	247021	33.458	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	255922	32.399	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	440074	30.120	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	561049	25.188	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	1421978	35.092	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1642860	35.564	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	43004	5.125	ng/u1	0.01
60) Fluorene-d10	15.522	176	1309917	37.223	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	202657	26.934	ng/u1	0.00
73) Anthracene-d10	17.387	188	2203715	38.917	ng/u1	0.00
81) Pyrene-d10	19.622	212	2801844	39.074	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.645	264	2951783	40.312	ng/u1	0.00

Target Compounds Qvalue

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

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