

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012393.D
 Acq On : 31 Oct 2022 18:58
 Operator : CG/JU
 Sample : N5194-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB900

Quant Time: Oct 31 23:52:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 28 06:05:59 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	274247	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.599	136	1215409	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.446	164	778493	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.210	188	1762558	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	2068911	20.000	ng/ul	0.00
88) Perylene-d12	23.692	264	2168482	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	24029	3.528	ng/uL	0.00
4) Pyridine-d5	3.676	84	85092	4.349	ng/ul	0.00
7) Phenol-d5	6.969	99	109403	4.448	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	379803	25.865	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	352513	19.625	ng/ul	0.00
15) 4-Methylphenol-d8	8.511	113	218996	11.257	ng/ul	-0.01
21) Nitrobenzene-d5	8.963	128	242777	30.803	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.687	143	241079	29.469	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	407801	23.002	ng/ul	0.00
31) 4-Chloroaniline-d4	10.746	131	378376	14.334	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1592561	29.884	ng/ul	0.00
49) Acenaphthylene-d8	14.140	160	1766175	29.001	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	33765	3.348	ng/ul	0.00
60) Fluorene-d10	15.445	176	1453484	31.856	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	216652	23.609	ng/ul	-0.01
73) Anthracene-d10	17.310	188	2416764	31.823	ng/ul	0.00
81) Pyrene-d10	19.551	212	3091072	29.758	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.539	264	3355549	31.670	ng/ul	-0.01

Target Compounds Qvalue

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

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