

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012394.D
 Acq On : 31 Oct 2022 19:33
 Operator : CG/JU
 Sample : N5194-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB901

Quant Time: Oct 31 23:52:13 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	313196	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.598	136	1381501	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	895265	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.210	188	1982484	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	2268885	20.000	ng/ul	0.00
88) Perylene-d12	23.692	264	2276699	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	28370	3.647	ng/uL	0.00
4) Pyridine-d5	3.669	84	134260	6.009	ng/ul	0.00
7) Phenol-d5	6.969	99	129027	4.593	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	462254	27.565	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	417781	20.366	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	261474	11.769	ng/ul	-0.01
21) Nitrobenzene-d5	8.963	128	299356	33.415	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.687	143	289855	31.172	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	496502	24.638	ng/ul	0.00
31) 4-Chloroaniline-d4	10.745	131	634500	21.146	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1908086	31.134	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	2159015	30.827	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	36716	3.165	ng/ul	0.00
60) Fluorene-d10	15.445	176	1696188	32.326	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	247659	23.994	ng/ul	0.00
73) Anthracene-d10	17.304	188	2717723	31.816	ng/ul	-0.01
81) Pyrene-d10	19.551	212	3394161	29.796	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.539	264	3510879	31.561	ng/ul	-0.01

Target Compounds Qvalue

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

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