

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012434.D
 Acq On : 01 Nov 2022 19:01
 Operator : CG/JU
 Sample : N5270-02
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 01 23:00:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	255153	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	1179796	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	775935	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1690476	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	1654562	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	1449207	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	30207	4.767	ng/uL	0.00
4) Pyridine-d5	3.664	84	145203	7.977	ng/ul	0.00
7) Phenol-d5	6.963	99	211627	9.247	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	610279	44.671	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	593547	35.516	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	424190	23.436	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	413293	54.020	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	436062	54.913	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	747410	43.430	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	810197	31.618	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	2685937	50.567	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	3018113	49.721	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	72662	7.228	ng/ul	0.00
60) Fluorene-d10	15.440	176	2310497	50.805	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	404666	45.978	ng/ul	0.00
73) Anthracene-d10	17.304	188	3747576	51.450	ng/ul	0.00
81) Pyrene-d10	19.545	212	4387338	52.815	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.527	264	3801235	53.682	ng/ul	-0.02

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

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