

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012386.D
 Acq On : 31 Oct 2022 14:22
 Operator : CG/JU
 Sample : N5180-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0KX3

Quant Time: Oct 31 23:50:21 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	7.799	152	318171	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	1384814	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	884570	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	1952108	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	2166147	20.000	ng/ul	0.00
88) Perylene-d12	23.698	264	2209197	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	32170	4.071	ng/uL	0.00
4) Pyridine-d5	3.669	84	177197	7.806	ng/ul	0.00
7) Phenol-d5	6.975	99	71379	2.501	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	489403	28.728	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	248309	11.915	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	181247	8.030	ng/ul	0.00
21) Nitrobenzene-d5	8.969	128	311992	34.742	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	3646	0.391	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	233346	11.552	ng/ul	0.00
31) 4-Chloroaniline-d4	10.746	131	667756	22.202	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1968588	32.510	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	2240509	32.377	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	51	0.004	ng/ul	0.00
60) Fluorene-d10	15.445	176	1818274	35.072	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	4	0.000	ng/ul	0.00
73) Anthracene-d10	17.310	188	3277618	38.967	ng/ul	0.00
81) Pyrene-d10	19.551	212	3915617	36.004	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.545	264	4162647	38.563	ng/ul	0.00

Target Compounds Qvalue

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

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