

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012406.D
 Acq On : 01 Nov 2022 02:25
 Operator : CG/JU
 Sample : N5216-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA57

Quant Time: Nov 01 03:18:13 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	7.793	152	288500	20.000	ng/u1	0.00
20) Naphthalene-d8	10.593	136	1295469	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.446	164	842845	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	1865867	20.000	ng/u1	0.00
79) Chrysene-d12	21.298	240	2055948	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	2081359	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	27362	3.819	ng/uL	0.00
4) Pyridine-d5	3.670	84	99200	4.820	ng/u1	0.01
7) Phenol-d5	6.969	99	138003	5.333	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	443243	28.694	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	438028	23.180	ng/u1	0.00
15) 4-Methylphenol-d8	8.511	113	279400	13.652	ng/u1	0.00
21) Nitrobenzene-d5	8.963	128	294679	35.077	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	310803	35.645	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	515536	27.282	ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	718304	25.529	ng/u1	0.00
46) Dimethylphthalate-d6	13.863	166	2064767	35.786	ng/u1	0.00
49) Acenaphthylene-d8	14.140	160	2237010	33.927	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	59037	5.406	ng/u1	0.00
60) Fluorene-d10	15.440	176	1828769	37.020	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	340948	35.097	ng/u1	0.00
73) Anthracene-d10	17.304	188	3234532	40.232	ng/u1	0.00
81) Pyrene-d10	19.545	212	3888291	37.669	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	4121127	40.523	ng/u1	-0.01

Target Compounds Qvalue

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

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