

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012412.D
 Acq On : 01 Nov 2022 06:24
 Operator : CG/JU
 Sample : N5216-12
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHAC0

Quant Time: Nov 01 07:25:39 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	356671	20.000 ng/u1	0.00
20) Naphthalene-d8	10.593	136	1618733	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.445	164	1048732	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.204	188	2293720	20.000 ng/u1	0.00
79) Chrysene-d12	21.298	240	2530684	20.000 ng/u1	0.00
88) Perylene-d12	23.686	264	2525315	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.240	96	29386	3.317 ng/uL	0.00
4) Pyridine-d5	3.670	84	116784	4.590 ng/u1	0.01
7) Phenol-d5	6.964	99	184926	5.781 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	540005	28.276 ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	547825	23.450 ng/u1	0.00
15) 4-Methylphenol-d8	8.505	113	377117	14.905 ng/u1	0.00
21) Nitrobenzene-d5	8.963	128	370933	35.337 ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	393424	36.109 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	709356	30.042 ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	783345	22.281 ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2532284	35.273 ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	2822270	34.400 ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	88133	6.486 ng/u1	0.00
60) Fluorene-d10	15.440	176	2207785	35.919 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	413704	34.643 ng/u1	0.00
73) Anthracene-d10	17.304	188	3688109	37.317 ng/u1	0.00
81) Pyrene-d10	19.545	212	4454617	35.060 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	4669553	37.844 ng/u1	-0.01

Target Compounds Qvalue

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

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