

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012399.D
 Acq On : 31 Oct 2022 22:25
 Operator : CG/JU
 Sample : N5216-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA86

Quant Time: Oct 31 23:53:04 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	324811	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.592	136	1442323	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.445	164	925548	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.204	188	2045283	20.000	ng/ul	-0.01
79) Chrysene-d12	21.298	240	2249967	20.000	ng/ul	-0.01
88) Perylene-d12	23.686	264	2333936	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	29294	3.631	ng/uL	0.00
4) Pyridine-d5	3.669	84	111793	4.824	ng/ul	0.00
7) Phenol-d5	6.969	99	137316	4.713	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	464541	26.711	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	441289	20.742	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	267614	11.614	ng/ul	-0.01
21) Nitrobenzene-d5	8.963	128	306952	32.818	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.687	143	317448	32.700	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	520406	24.735	ng/ul	0.00
31) 4-Chloroaniline-d4	10.745	131	712113	22.732	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1995062	31.488	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	2210666	30.532	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	47628	3.972	ng/ul	0.00
60) Fluorene-d10	15.445	176	1787670	32.955	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	331507	31.131	ng/ul	-0.01
73) Anthracene-d10	17.304	188	3155488	35.806	ng/ul	-0.01
81) Pyrene-d10	19.545	212	3852680	34.106	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.539	264	4148733	36.380	ng/ul	-0.01

Target Compounds Qvalue

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

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