

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012136.D
 Acq On : 14 Oct 2022 14:07
 Operator : CG/JU
 Sample : N4982-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 14 23:03:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	263267	20.000	ng/ul	0.00
20) Naphthalene-d8	10.652	136	1115661	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.493	164	644776	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	1353876	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1379920	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	1465057	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	23937	3.585	ng/uL	0.00
4) Pyridine-d5	3.705	84	113111	6.370	ng/ul	0.00
7) Phenol-d5	7.016	99	156382	7.386	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	396058	32.954	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	442002	26.299	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	268964	16.538	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128	274092	34.786	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143	271992	34.044	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	435038	28.096	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	478496	20.970	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1344981	32.937	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	1642626	32.530	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	34283	4.276	ng/ul	0.00
60) Fluorene-d10	15.492	176	1245880	33.722	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	168155	23.522	ng/ul	0.00
73) Anthracene-d10	17.351	188	2019464	34.350	ng/ul	0.00
81) Pyrene-d10	19.592	212	2390455	33.814	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	2461822	34.567	ng/ul	0.00

Target Compounds Qvalue

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

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