

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021622\
 Data File : BP009178.D
 Acq On : 16 Feb 2022 14:57
 Operator : CG/JU
 Sample : N1449-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SPLP-LEACHATE-AMENDED-COMP

Quant Time: Feb 17 01:57:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 16 14:30:24 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 02/17/2022
 Supervised By :Jagrut Upadhyay 02/17/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	175022	20.000	ng	0.00
21) Naphthalene-d8	10.581	136	636356	20.000	ng	0.00
39) Acenaphthene-d10	14.428	164	421939	20.000	ng	0.00
64) Phenanthrene-d10	17.186	188	867791	20.000	ng	0.00
76) Chrysene-d12	21.286	240	963054	20.000	ng	0.00
86) Perylene-d12	23.674	264	1119762	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	613059	62.482	ng	0.00
7) Phenol-d6	6.981	99	414427	32.055	ng	0.00
23) Nitrobenzene-d5	8.946	82	1050046	93.055	ng	0.00
42) 2,4,6-Tribromophenol	15.934	330	932053	165.444	ng	0.00
45) 2-Fluorobiphenyl	13.051	172	2906480	96.430	ng	0.00
79) Terphenyl-d14	19.751	244	4370833	81.617	ng	0.00
Target Compounds						
32) Benzoic acid	9.981	122	9175m	4.241	ng	Qvalue

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

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